Example #	R_1	Name	NMR Data*	Mass Spec.
48		2,2'-[[4-(4-Morpholinyl)phenyl]methylene]bis[4-[[(5-methyl-1H-tetrazol-1-yl)imino]methyl]]phenol	^{13}H NMR in DMSO: 10.47(s, 2H); 9.15(s, 2H); 7.78(dd, J=1.76, 2.34, 8.50Hz, 2H); 7.46 (d, J=1.76Hz, 2H); 6.94 (m, 6H); 5.99 (s, 1H); 3.73 (t, J=4.69Hz, 4H); 3.07 (t, J=4.69Hz, 4H); 2.47 (s, 6H)	(M+H)+ 580.1
49		4-[[Bis[[5-(5-methyl-1H-tetrazol-1-yl)imino]methyl]-2-hydroxyphenyl]methylene]benzeneethanol, acetate ester	^{1}H NMR in DMSO: 520 (s. 2H); 9.15 (s, 2H); 7.80 (dd, J = 1.7, 8.2, 8.8Hz, 2H); 7.45 (d, J = 1.7Hz, 2H); 1 (d, J = 8.2Hz, 2H); 7.03 (d, J = 8.2Hz, 2H); 9 (d, J = 8.2Hz, 2H); 6.05 (s, 1H); 4.20 (t, J = Hz, 2H); 2.87 (t, J = 6.4, 7.0Hz, 2H); 2.46 (s,); 1.98 (s, 3H)	580, 552, 525, 456
50		2,2'-[[4-(1-Piperidinyl)phenyl]methylene]bis[4-[[(5-methyl-1H-tetrazol-1-yl)imino]methyl]]phenol	^{1}H NMR in DMSO:) 10.48 (s, 2H); 9.15 (s, 2H); 7.80 (d, J=8.21 Hz, 2H); 7.47 (s, 2H); 7.00 (d, J=8.79 Hz, 2H); 6.93 (s, 2H); 5.98 (s, 1H); 3.11 (m, 4H); 2.47 (s. 6H); 1.62 (m, 4H); 1.54 m. 2H)	576 (M-H)

*All ^{1}H NMR and ^{13}C NMR spectra were acquired on a Varian Mercury VX 300 Spectrometer and referenced to tetramethylsilane (TMS) unless indicated otherwise. Chemical shifts and coupling constants are reported in parts per million (ppm) and Hertz (Hz) respectively.